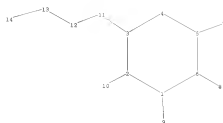
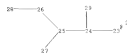
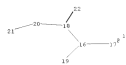
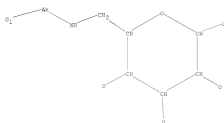
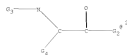
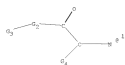


C

38



```

chain nodes :
  8 9 11 12 13 14 16 17 18 19 20 21 22 23 24 25 26 27 28 29 38
ring nodes :
  1 2 3 4 5 6
ring/chain nodes :
  7 10
chain bonds :
  1-9 3-11 6-8 11-12 12-13 13-14 16-17 16-18 16-19 18-20 18-22 20-21 23-24 24-25
24-29 25-26 25-27 26-28
ring/chain bonds :
  2-10 5-7
ring bonds :
  1-2 1-6 2-3 3-4 4-5 5-6
exact/norm bonds :
  1-2 1-6 1-9 2-3 2-10 3-4 4-5 5-6 5-7 6-8 12-13 13-14 16-17 16-19 18-20 18-22
20-21 23-24 24-29 25-26 25-27 26-28
exact bonds :
  3-11 11-12 16-18 24-25

```

G1:[\*1],[\*2]

G2:O,S,N

G3:H,Cy,Ak

G4:H,CH

```

Match level :
  1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11:CLASS
12:CLASS 13:CLASS 14:CLASS 16:CLASS 17:CLASS 18:CLASS 19:CLASS 20:CLASS 21:CLASS
22:CLASS 23:CLASS 24:CLASS 25:CLASS 26:CLASS 27:CLASS 28:CLASS 29:CLASS 38:CLASS

```

Element Count :

```

Node 13: Limited
C,C1-20

```

